

# STN SEARCH TRANSCRIPT

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\* \* \* \* \* \* \* \* \* \* \* Welcome to STN International \* \* \* \* \* \* \* \* \* \* \*

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NEWS 8 DEC 23 New IPC8 SEARCH, DISPLAY, and SELECT fields in USPATFULL/  
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NEWS 9 JAN 13 IPC 8 searching in IFIPAT, IFIUDB, and IFICDB  
NEWS 10 JAN 13 New IPC 8 SEARCH, DISPLAY, and SELECT enhancements added to  
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NEWS 11 JAN 17 Pre-1988 INPI data added to MARPAT  
NEWS 12 JAN 17 IPC 8 in the WPI family of databases including WPIFV  
NEWS 13 JAN 30 Saved answer limit increased  
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added to TULSA

NEWS EXPRESS JANUARY 03 CURRENT VERSION FOR WINDOWS IS V8.01,  
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=> FILE REG  
COST IN U.S. DOLLARS  
  
FULL ESTIMATED COST

SINCE FILE ENTRY	0.21	TOTAL SESSION	0.21
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FILE 'REGISTRY' ENTERED AT 10:04:57 ON 08 FEB 2006  
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STRUCTURE FILE UPDATES: 6 FEB 2006 HIGHEST RN 873652-66-5  
DICTIONARY FILE UPDATES: 6 FEB 2006 HIGHEST RN 873652-66-5

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TSCA INFORMATION NOW CURRENT THROUGH JULY 14, 2005

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\*\*\*\*\*  
\*  
\* The CA roles and document type information have been removed from \*  
\* the IDE default display format and the ED field has been added, \*  
\* effective March 20, 2005. A new display format, IDERL, is now \*  
\* available and contains the CA role and document type information. \*  
\*  
\*\*\*\*\*

Structure search iteration limits have been increased. See HELP SLIMITS  
for details.

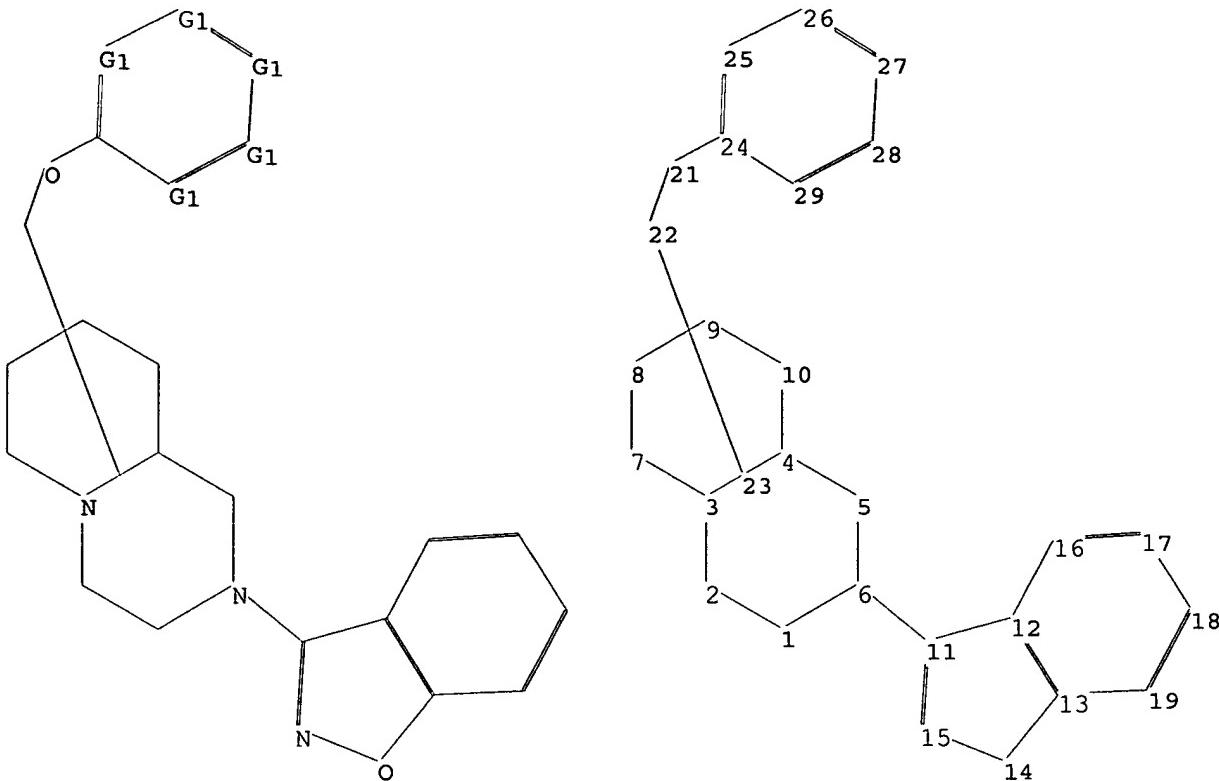
REGISTRY includes numerically searchable data for experimental and  
predicted properties as well as tags indicating availability of  
experimental property data in the original document. For information  
on property searching in REGISTRY, refer to:

<http://www.cas.org/ONLINE/UG/regprops.html>

=> ....Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END) :end

=>  
Uploading C:\Program Files\Stnexp\Queries\BRIGHT ANTISSYCHOTICS.str



chain nodes :

21 22

ring nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 24 25 26 27  
28 29

chain bonds :

6-11 21-22 21-24

ring bonds :

1-2 1-6 2-3 3-4 3-7 4-5 4-10 5-6 7-8 8-9 9-10 11-12 11-15 12-13 12-16  
13-14 13-19 14-15 16-17 17-18 18-19 24-25 24-29 25-26 26-27 27-28 28-29

exact/norm bonds :

1-2 1-6 2-3 3-4 3-7 4-5 4-10 5-6 6-11 7-8 8-9 9-10 11-12 11-15 13-14  
14-15 21-22 21-24 24-25 24-29 25-26 26-27 27-28 28-29

normalized bonds :

12-13 12-16 13-19 16-17 17-18 18-19

isolated ring systems :

containing 1 : 11 : 24 :

G1:C,N

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom

11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom

21:CLASS 22:CLASS

23:CLASS 24:Atom 25:Atom 26:Atom 27:Atom 28:Atom 29:Atom

L1 STRUCTURE UPLOADED

=> que L1

L2 QUE L1

=> D L1  
L1 HAS NO ANSWERS  
L1 STR

\* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT \*  
Structure attributes must be viewed using STN Express query preparation.

=> S L1  
SAMPLE SEARCH INITIATED 10:05:18 FILE 'REGISTRY'  
SAMPLE SCREEN SEARCH COMPLETED - 13 TO ITERATE

100.0% PROCESSED 13 ITERATIONS 13 ANSWERS  
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*  
BATCH \*\*COMPLETE\*\*  
PROJECTED ITERATIONS: 44 TO 476  
PROJECTED ANSWERS: 44 TO 476

L3 13 SEA SSS SAM L1

=> S L1 SSS FULL  
FULL SEARCH INITIATED 10:05:23 FILE 'REGISTRY'  
FULL SCREEN SEARCH COMPLETED - 171 TO ITERATE

100.0% PROCESSED 171 ITERATIONS 168 ANSWERS  
SEARCH TIME: 00.00.01

L4 168 SEA SSS FUL L1

=> FILE CAPLUS  
COST IN U.S. DOLLARS SINCE FILE TOTAL  
ENTRY SESSION  
FULL ESTIMATED COST 166.94 167.15

FILE 'CAPLUS' ENTERED AT 10:05:26 ON 08 FEB 2006  
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FILE COVERS 1907 - 8 Feb 2006 VOL 144 ISS 7  
FILE LAST UPDATED: 7 Feb 2006 (20060207/ED)

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=> S L4  
L5 3 L4

=> D 1-3 IBIB ABS

L5 ANSWER 1 OF 3 CAPLUS COPYRIGHT 2006 ACS on STN  
ACCESSION NUMBER: 2005:698356 CAPLUS  
DOCUMENT NUMBER: 143:179645  
TITLE: Compositions containing atypical antipsychotics and azabicyclic compounds for treating CNS disorders  
INVENTOR(S): Brodney, Michael A.; Howard, Harry R.  
PATENT ASSIGNEE(S): Pfizer Inc, USA  
SOURCE: U.S. Pat. Appl. Publ., 21 pp.  
CODEN: USXXCO  
DOCUMENT TYPE: Patent  
LANGUAGE: English  
FAMILY ACC. NUM. COUNT: 1  
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
US 2005171086	A1	20050804	US 2005-48013	20050128
WO 2005082370	A1	20050909	WO 2005-IB106	20050117
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SM, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW RW: BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IS, IT, LT, LU, MC, NL, PL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				

PRIORITY APPLN. INFO.: US 2004-539939P P 20040129  
OTHER SOURCE(S): MARPAT 143:179645  
AB Disclosed is an aminomethylpyridyloxymethyl/benzisoxazole substituted azabicyclic compound, a pharmaceutical composition comprising same, and a method of treating one or more CNS or other disorders, including concurrent treatment of disorders such as schizophrenia and depression. For example, capsules for Parkinson's disease contained ziprasidone hydrochloride 200, benzisoxazole substituted azabicyclic compd 20, Methocel E3 222, lactose monohydrate 222, Aerosil 10, SLS 10 mg.

L5 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2006 ACS on STN  
ACCESSION NUMBER: 2004:780698 CAPLUS  
DOCUMENT NUMBER: 141:296048  
TITLE: A preparation of pyrido[1,2-a]pyrazine derivatives, useful for the treatment of schizophrenia and depression  
INVENTOR(S): Bright, Gene Michael; Brodney, Michael Aaron;  
Wlodecki, Bishop  
PATENT ASSIGNEE(S): Pfizer Products Inc., USA  
SOURCE: PCT Int. Appl., 87 pp.  
CODEN: PIXXD2  
DOCUMENT TYPE: Patent  
LANGUAGE: English  
FAMILY ACC. NUM. COUNT: 1  
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2004081007	A1	20040923	WO 2004-IB499	20040223
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD,				

GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC,	
LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NA, NI,	
NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY,	
TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW	
RW: BW, GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ,	
BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE,	
ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK,	
TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG	
CA 2518740 AA 20040923 CA 2004-2518740 20040223	
EP 1608648 A1 20051228 EP 2004-713592 20040223	
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT,	
IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, SK	
NL 1025710 A1 20040914 NL 2004-1025710 20040312	
NL 1025710 C2 20051010	
US 2005026922 A1 20050203 US 2004-800328 20040312	
NO 2005004095 A 20050926 NO 2005-4095 20050902	
PRIORITY APPLN. INFO.: US 2003-453925P P 20030312	
	WO 2004-IB499 W 20040223

OTHER SOURCE(S) : MARPAT 141:296048  
GI

\* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT \*

AB The invention relates to a preparation of pyrido[1,2-a]pyrazine derivs. of formula I [wherein: X is O, NH, or N(alkyl); Y is (CH<sub>2</sub>)<sub>0-1</sub>; Z is CHO, C(O)-alkoxy, SO<sub>2</sub>-alkoxy, Me, CH<sub>2</sub>OH, etc.; R<sub>1</sub> and R<sub>2</sub> are independently selected from H, halogen, (cyclo)alkyl, or alkoxy, etc.], useful for treating CNS or other disorders, including concurrent treatment of disorders such as schizophrenia and depression. Thus, e.g., II was prepared via reaction of morpholine with (7R,9aS)-trans-methanesulfonic acid 6-(2-benzo[d]isoxazol-3-yl-octahydropyrido[1,2-a]pyrazin-7-ylmethoxy)-pyridin-2-ylmethyl ester (preparation given). The prepared compds. were determined to

be antagonists and/or inverse agonists of human D<sub>2</sub>, human 5-HT1B, and human 5-HT2A receptors. For instance, preferred compound II exhibited Ki value of about 20 nM or less for at least two of the following receptors: D<sub>2</sub>, 5-HT1B, and 5-HT2A.

REFERENCE COUNT: 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L5 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2006 ACS on STN  
 ACCESSION NUMBER: 1999:672814 CAPLUS  
 DOCUMENT NUMBER: 131:299376  
 TITLE: Azabicyclic 5-HT1 receptor ligands, particularly 2-(benzo[d]isoxazol-3-yl)-7-(phenoxy)methyl octahydropyrido[1,2-a]pyrazine derivatives  
 INVENTOR(S) : Bright, Gene Michael  
 PATENT ASSIGNEE(S) : Pfizer Products Inc., USA  
 SOURCE: PCT Int. Appl., 96 pp.  
 CODEN: PIXXD2  
 DOCUMENT TYPE: Patent  
 LANGUAGE: English  
 FAMILY ACC. NUM. COUNT: 1  
 PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
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WO 9952907	A1	19991021	WO 1999-IB457	19990318
W: AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CU, CZ, DE, DK, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IS, JP, KE,				

KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, UA, UG, US, UZ, VN, YU, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM RW: GH, GM, KE, LS, MW, SD, SL, SZ, UG, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG				
CA 2327782	AA	19991021	CA 1999-2327782	19990318
AU 9932687	A1	19991101	AU 1999-32687	19990318
AU 749254	B2	20020620		
BR 9909522	A	20001219	BR 1999-9522	19990318
TR 200002932	T2	20010122	TR 2000-200002932	19990318
EP 1070065	A1	20010124	EP 1999-945686	19990318
EP 1070065	B1	20030903		
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JP 2002511469	T2	20020416	JP 2000-543464	19990318
JP 3496759	B2	20040216		
NZ 507010	A	20030530	NZ 1999-507010	19990318
CN 1117092	B	20030806	CN 1999-807195	19990318
AT 248838	E	20030915	AT 1999-945686	19990318
PT 1070065	T	20031231	PT 1999-945686	19990318
ES 2204154	T3	20040416	ES 1999-945686	19990318
TW 555758	B	20031001	TW 1999-88105301	19990402
ZA 9902596	A	20001008	ZA 1999-2596	19990408
US 6525048	B1	20030225	US 2000-403892	20000118
NO 2000005004	A	20001127	NO 2000-5004	20001004
HR 2000000653	A1	20010430	HR 2000-653	20001004
BG 104915	A	20010731	BG 2000-104915	20001107
HK 1035719	A1	20031224	HK 2001-106215	20010904
US 2003181458	A1	20030925	US 2003-364804	20030211
US 6887905	B2	20050503		
PRIORITY APPLN. INFO.:			US 1998-81237P	P 19980409
			WO 1999-IB457	W 19990318
			US 2000-403892	A3 20000118

OTHER SOURCE(S) : MARPAT 131:299376  
GI

\* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT \*

AB The invention relates to compds. I [wherein R3, R4, Z = H, (fluoro)alkyl, (fluoro)alkoxy, (fluoro)alkoxy(fluoro)alkyl; W = alkoxy(methyl, (di)(alkyl)aminomethyl, or CH<sub>2</sub>NR1R2 where R1R2 = atoms to complete a heterocycle such as pyrrolidine; with provisos]. The compds. are agonists or antagonists of serotonin 1A receptors, and/or antagonists of serotonin 1D receptors, and are thus useful as psychotherapeutic agents. These compds. may be co-administered with 5-HT reuptake inhibitors, and are potentially useful for treating a wide variety of conditions. Approx. 40 synthetic examples are given. For instance, title compound II was prepared in 5 steps: (1) Mitsunobu etherification of starting material (7R,9aS)-trans-III with Me 3-hydroxybenzoate (75%); (2) reduction of the Me ester to an alc. using LiAlH<sub>4</sub> (100%); (3) mesylation of the alc. and reaction of the mesylate with pyrrolidine (56%); (4) removal of the BOC protecting group (100%); and coupling with 3-chloro-5-fluorobenzo[d]isoxazole (36%). In assays against 5-HT receptors in vitro, all tested compds. I exhibited IC<sub>50</sub> values of < 0.60 mM for 5-HT1D receptors, and < 1.0 mM for 5-HT1A receptors.

REFERENCE COUNT: 1 THERE ARE 1 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

=> D 3 HITSTR

L5 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2006 ACS on STN

IT 247091-72-1P 247091-73-2P 247091-74-3P  
247091-75-4P 247091-76-5P 247091-77-6P  
247091-78-7P 247091-79-8P 247091-80-1P  
247091-81-2P 247091-82-3P 247091-83-4P  
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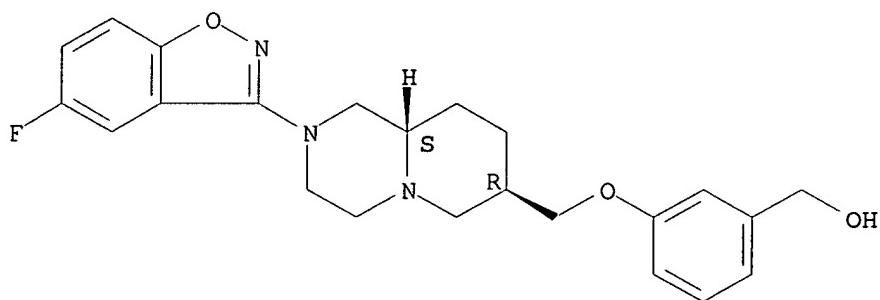
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(intermediate; preparation of benzisoxazolyloctahydropyridopyrazine derivs. as 5-HT1 receptor ligands)

RN 247091-72-1 CAPLUS

CN Benzenemethanol, 3-[[[(7R,9aS)-2-(5-fluoro-1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl)methoxy]- (9CI) (CA INDEX NAME)

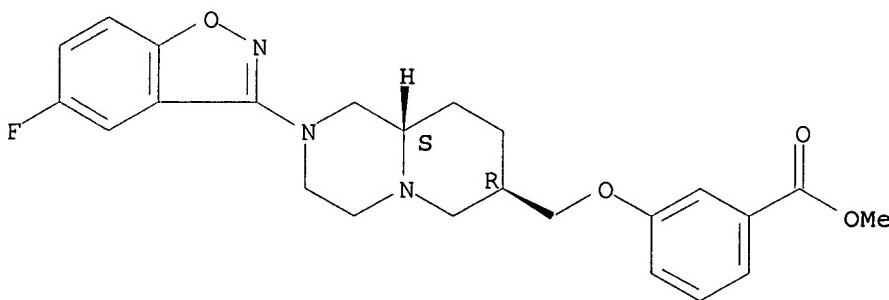
Absolute stereochemistry.



RN 247091-73-2 CAPLUS

CN Benzoic acid, 3-[[[(7R,9aS)-2-(5-fluoro-1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl)methoxy]-, methyl ester (9CI) (CA INDEX NAME)

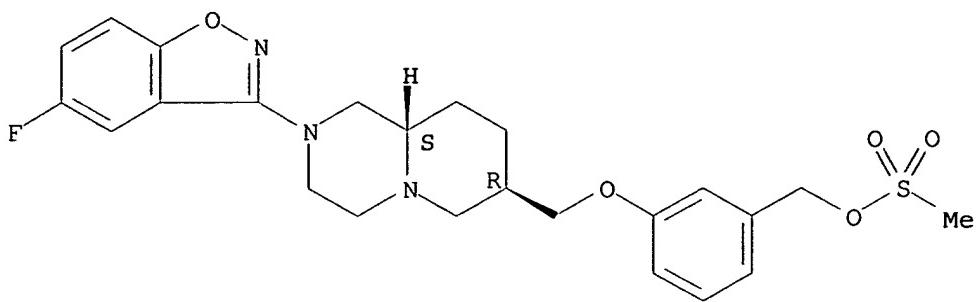
Absolute stereochemistry.



RN 247091-74-3 CAPLUS

CN Benzenemethanol, 3-[[[(7R,9aS)-2-(5-fluoro-1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl)methoxy]-, methanesulfonate (ester) (9CI) (CA INDEX NAME)

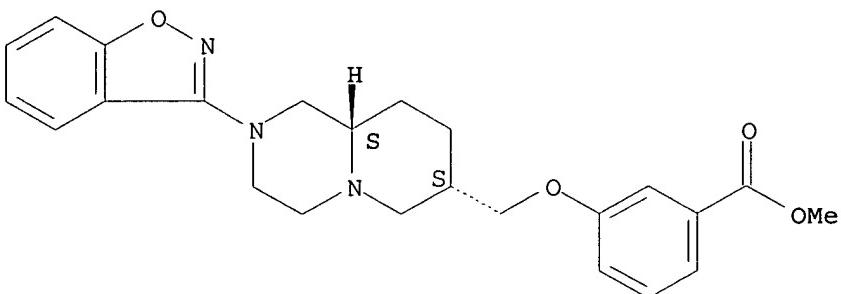
Absolute stereochemistry.



RN 247091-75-4 CAPLUS

CN Benzoic acid, 3-[(7S,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy]-, methyl ester (9CI) (CA INDEX NAME)

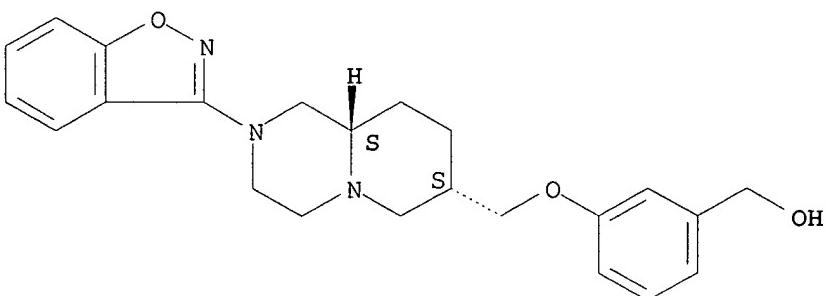
Absolute stereochemistry.



RN 247091-76-5 CAPLUS

CN Benzenemethanol, 3-[(7S,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy]- (9CI) (CA INDEX NAME)

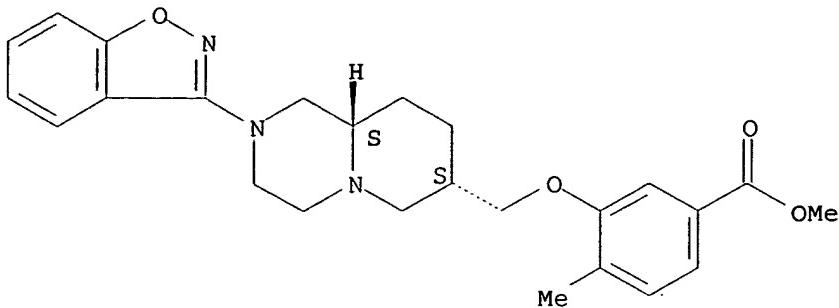
Absolute stereochemistry.



RN 247091-77-6 CAPLUS

CN Benzoic acid, 3-[(7S,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy]-4-methyl-, methyl ester (9CI) (CA INDEX NAME)

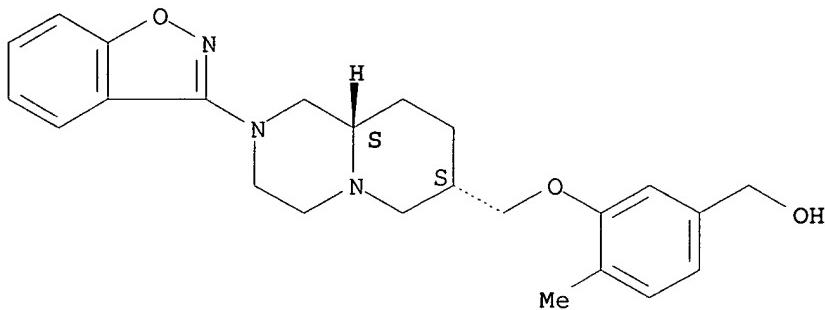
Absolute stereochemistry.



RN 247091-78-7 CAPLUS

CN Benzenemethanol, 3-[(7S,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy]-4-methyl- (9CI) (CA INDEX NAME)

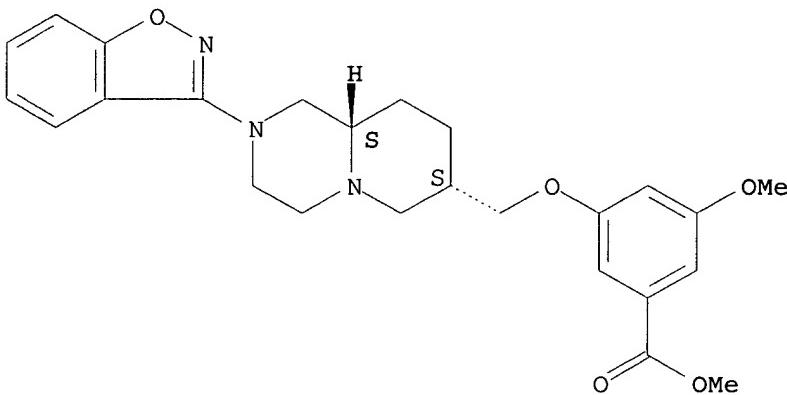
Absolute stereochemistry.



RN 247091-79-8 CAPLUS

CN Benzoic acid, 3-[(7S,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy]-5-methoxy-, methyl ester (9CI) (CA INDEX NAME)

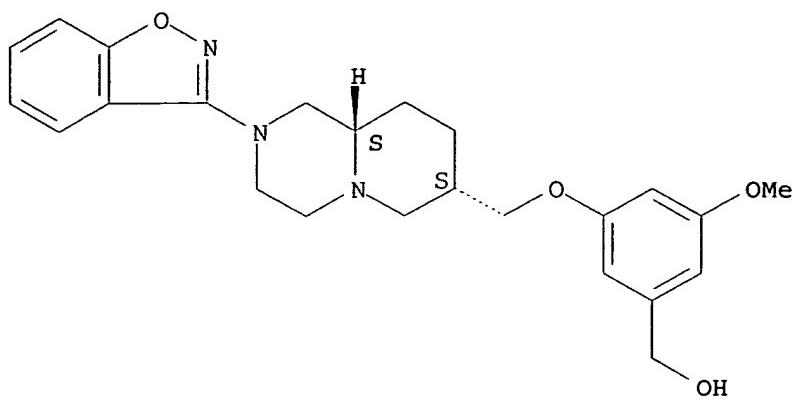
Absolute stereochemistry.



RN 247091-80-1 CAPLUS

CN Benzenemethanol, 3-[(7S,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy]-5-methoxy- (9CI) (CA INDEX NAME)

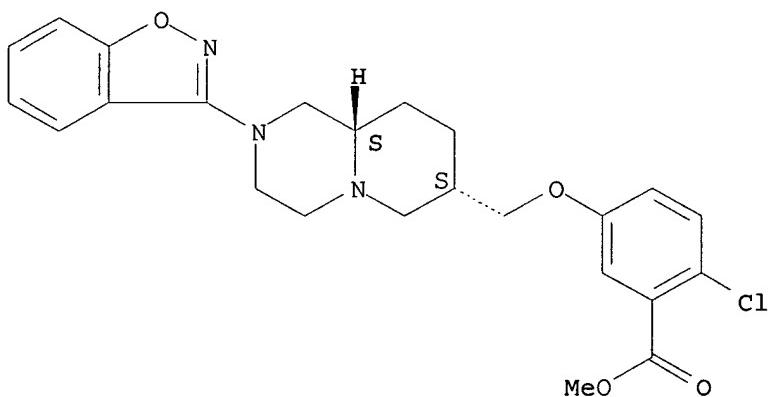
Absolute stereochemistry.



RN 247091-81-2 CAPLUS

CN Benzoic acid, 5-[(7S,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy]-2-chloro-, methyl ester (9CI) (CA INDEX NAME)

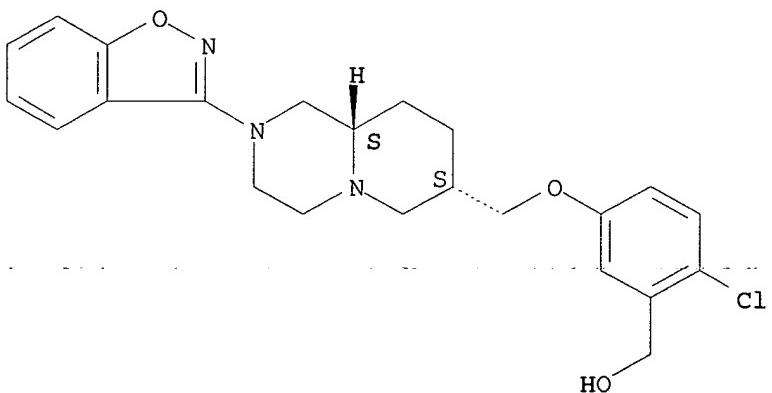
Absolute stereochemistry.



RN 247091-82-3 CAPLUS

CN Benzenemethanol, 5-[(7S,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy]-2-chloro- (9CI) (CA INDEX NAME)

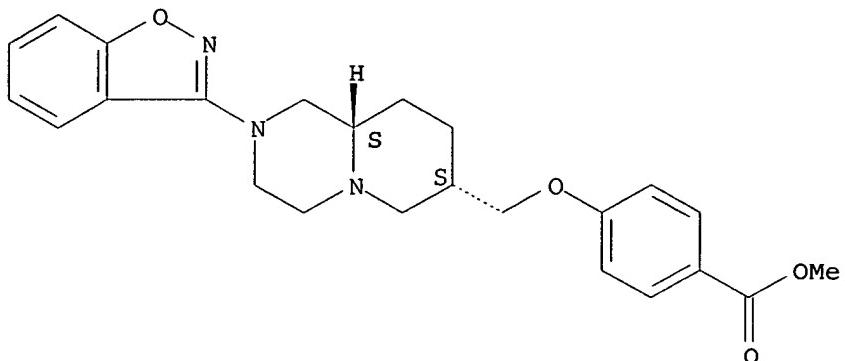
Absolute stereochemistry.



RN 247091-83-4 CAPLUS

CN Benzoic acid, 4-[(7S,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy-, methyl ester (9CI) (CA INDEX NAME)

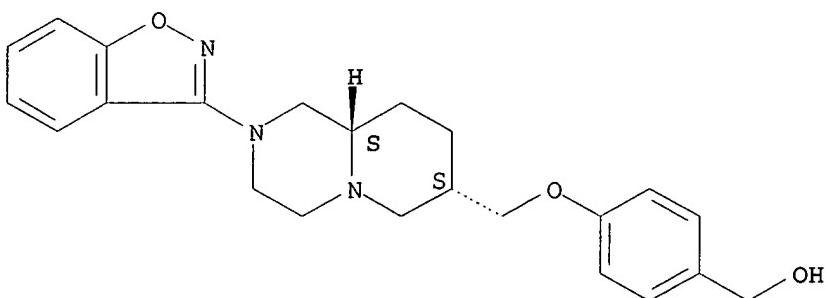
Absolute stereochemistry.



RN 247091-84-5 CAPLUS

CN Benzenemethanol, 4-[(7S,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy]- (9CI) (CA INDEX NAME)

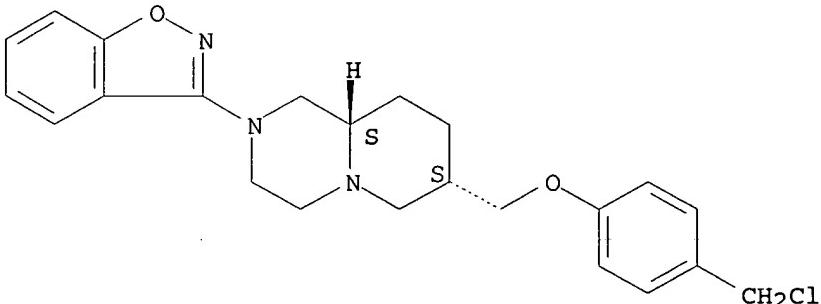
Absolute stereochemistry.



RN 247091-85-6 CAPLUS

CN 2H-Pyrido[1,2-a]pyrazine, 2-(1,2-benzisoxazol-3-yl)-7-[[4-(chloromethyl)phenoxy]methyl]octahydro-, (7S,9aS)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

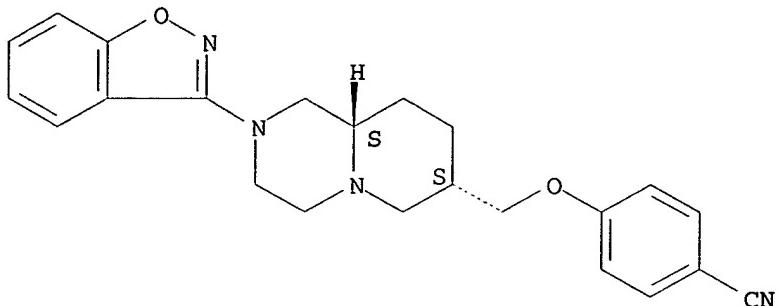


RN 247091-88-9 CAPLUS

CN Benzonitrile, 4-[(7S,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2H-

pyrido[1,2-a]pyrazin-7-ylmethoxy] - (9CI) (CA INDEX NAME)

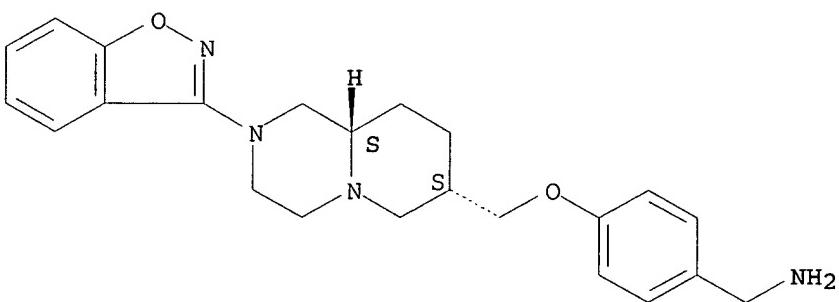
Absolute stereochemistry.



RN 247091-89-0 CAPLUS

CN Benzenemethanamine, 4-[[[(7S,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy] - (9CI) (CA INDEX NAME)

Absolute stereochemistry.



IT 247091-25-4P 247091-26-5P 247091-27-6P

247091-28-7P 247091-29-8P 247091-30-1P

247091-31-2P 247091-32-3P 247091-33-4P

247091-34-5P 247091-35-6P 247091-36-7P

247091-37-8P 247091-38-9P 247091-39-0P

247091-40-3P 247091-41-4P 247091-42-5P

247091-43-6P 247091-44-7P 247091-45-8P

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247091-50-5P 247091-51-6P 247091-52-7P

247091-53-8P

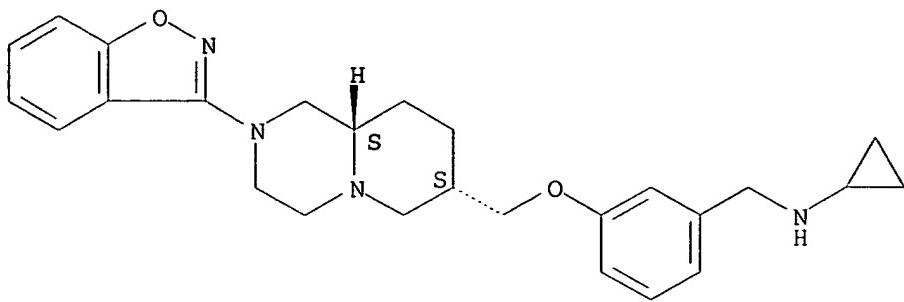
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)  
(target compound; preparation of benzisoxazolyloctahydropyridopyrazine derivs.

as 5-HT1 receptor ligands)

RN 247091-25-4 CAPLUS

CN Benzenemethanamine, 3-[[[(7S,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy]-N-cyclopropyl- (9CI) (CA INDEX NAME)

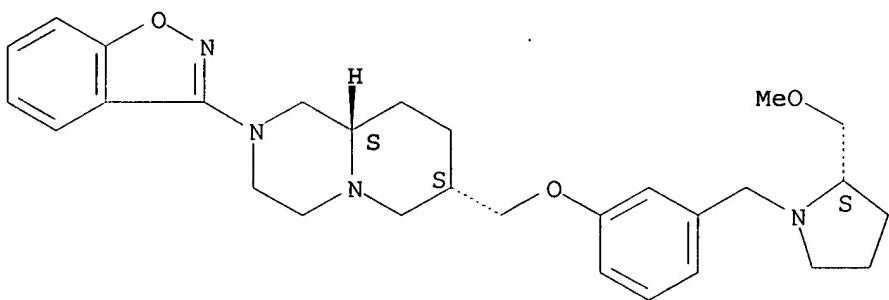
Absolute stereochemistry.



RN 247091-26-5 CAPLUS

CN 2H-Pyrido[1,2-a]pyrazine, 2-(1,2-benzisoxazol-3-yl)octahydro-7-[[3-[(2S)-2-(methoxymethyl)-1-pyrrolidinylmethyl]phenoxy]methyl]-, (7S,9aS)- (9CI) (CA INDEX NAME)

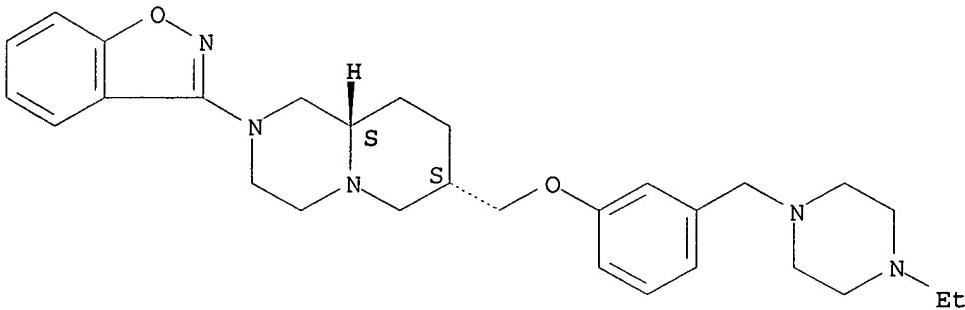
Absolute stereochemistry.



RN 247091-27-6 CAPLUS

CN 2H-Pyrido[1,2-a]pyrazine, 2-(1,2-benzisoxazol-3-yl)-7-[[3-[(4-ethyl-1-piperazinyl)methyl]phenoxy]methyl]octahydro-, (7S,9aS)- (9CI) (CA INDEX NAME)

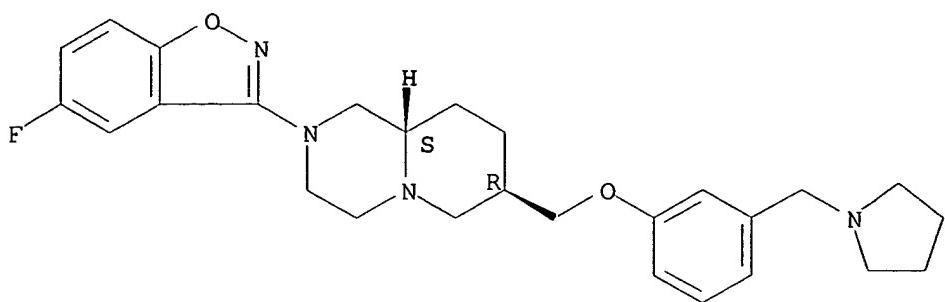
Absolute stereochemistry.



RN 247091-28-7 CAPLUS

CN 2H-Pyrido[1,2-a]pyrazine, 2-(5-fluoro-1,2-benzisoxazol-3-yl)octahydro-7-[[3-(1-pyrrolidinylmethyl)phenoxy]methyl]-, (7R,9aS)- (9CI) (CA INDEX NAME)

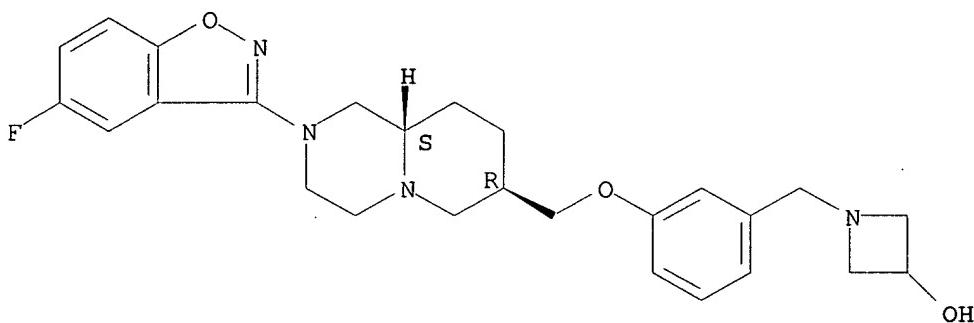
Absolute stereochemistry.



RN 247091-29-8 CAPLUS

CN 3-Azetidinol, 1-[[3-[(7R,9aS)-2-(5-fluoro-1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy]phenyl]methyl]- (9CI) (CA INDEX NAME)

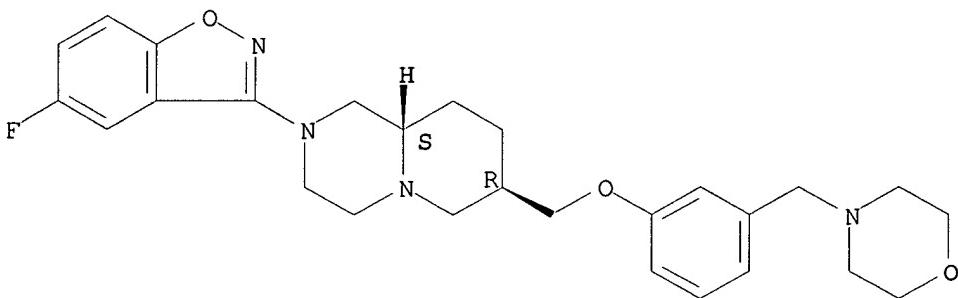
Absolute stereochemistry.



RN 247091-30-1 CAPLUS

CN 2H-Pyrido[1,2-a]pyrazine, 2-(5-fluoro-1,2-benzisoxazol-3-yl)octahydro-7-[[3-(4-morpholinylmethyl)phenoxy]methyl]-, (7R,9aS)- (9CI) (CA INDEX NAME)

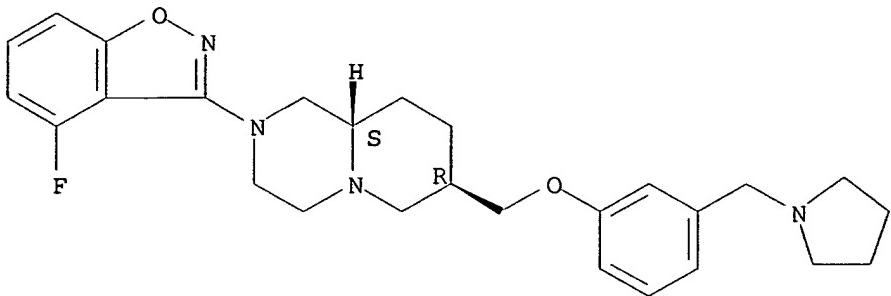
Absolute stereochemistry.



RN 247091-31-2 CAPLUS

CN 2H-Pyrido[1,2-a]pyrazine, 2-(4-fluoro-1,2-benzisoxazol-3-yl)octahydro-7-[[3-(1-pyrrolidinylmethyl)phenoxy]methyl]-, (7R,9aS)- (9CI) (CA INDEX NAME)

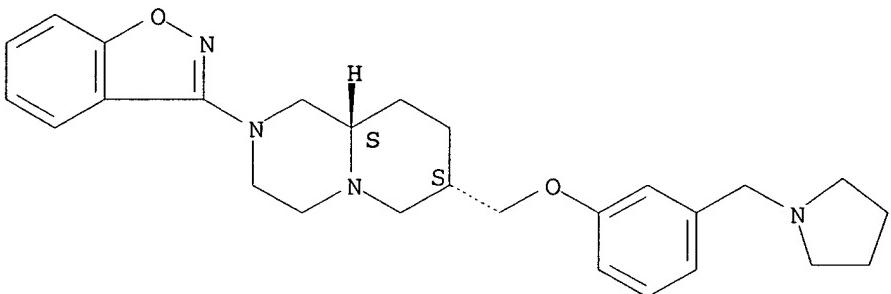
Absolute stereochemistry.



RN 247091-32-3 CAPLUS

CN 2H-Pyrido[1,2-a]pyrazine, 2-(1,2-benzisoxazol-3-yl)octahydro-7-[(3-(1-pyrrolidinylmethyl)phenoxy)methyl]-, (7S,9aS)- (9CI) (CA INDEX NAME)

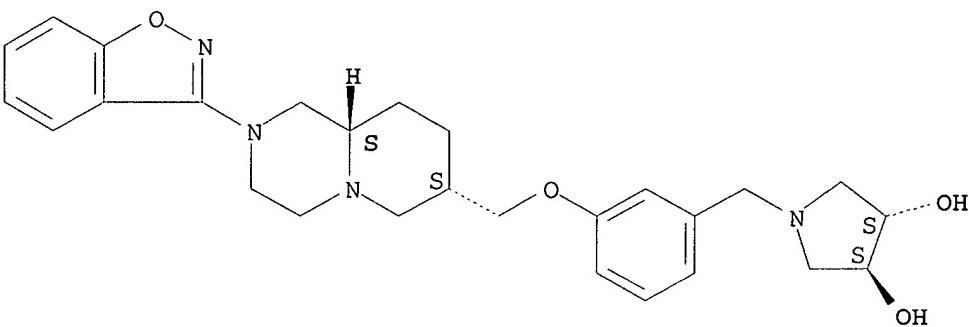
Absolute stereochemistry.



RN 247091-33-4 CAPLUS

CN 3,4-Pyrrolidinediol, 1-[[3-[(7S,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy]phenyl]methyl-, (3S,4S)- (9CI) (CA INDEX NAME)

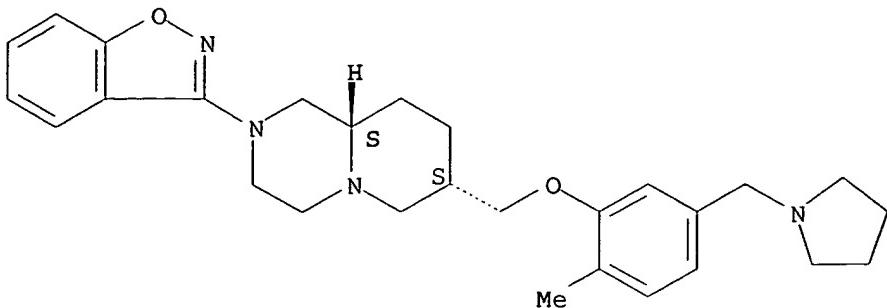
Absolute stereochemistry.



RN 247091-34-5 CAPLUS

CN 2H-Pyrido[1,2-a]pyrazine, 2-(1,2-benzisoxazol-3-yl)octahydro-7-[(2-methyl-5-(1-pyrrolidinylmethyl)phenoxy)methyl]-, (7S,9aS)- (9CI) (CA INDEX NAME)

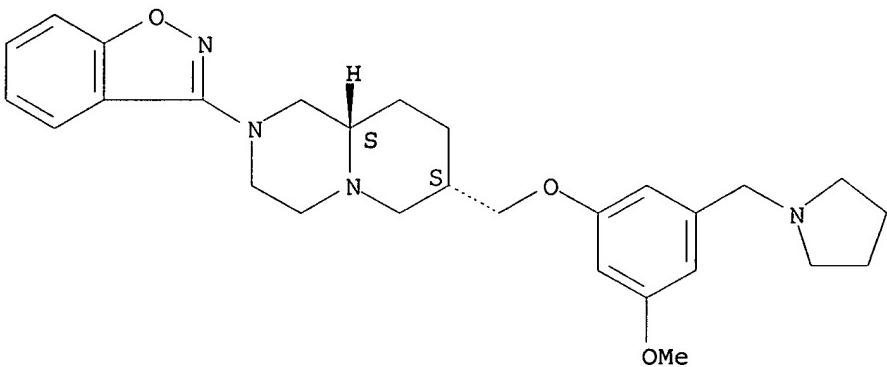
Absolute stereochemistry.



RN 247091-35-6 CAPLUS

CN 2H-Pyrido[1,2-a]pyrazine, 2-(1,2-benzisoxazol-3-yl)octahydro-7-[[3-methoxy-5-(1-pyrrolidinylmethyl)phenoxy]methyl]-, (7S,9aS)- (9CI) (CA INDEX NAME)

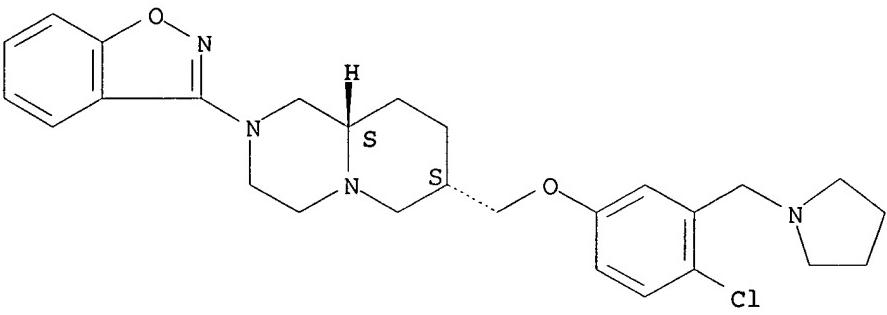
Absolute stereochemistry.



RN 247091-36-7 CAPLUS

CN 2H-Pyrido[1,2-a]pyrazine, 2-(1,2-benzisoxazol-3-yl)-7-[[4-chloro-3-(1-pyrrolidinylmethyl)phenoxy]methyl]octahydro-, (7S,9aS)- (9CI) (CA INDEX NAME)

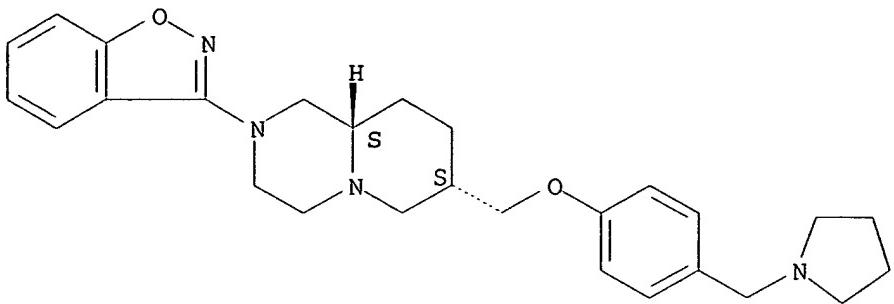
Absolute stereochemistry.



RN 247091-37-8 CAPLUS

CN 2H-Pyrido[1,2-a]pyrazine, 2-(1,2-benzisoxazol-3-yl)octahydro-7-[[4-(1-pyrrolidinylmethyl)phenoxy]methyl]-, (7S,9aS)- (9CI) (CA INDEX NAME)

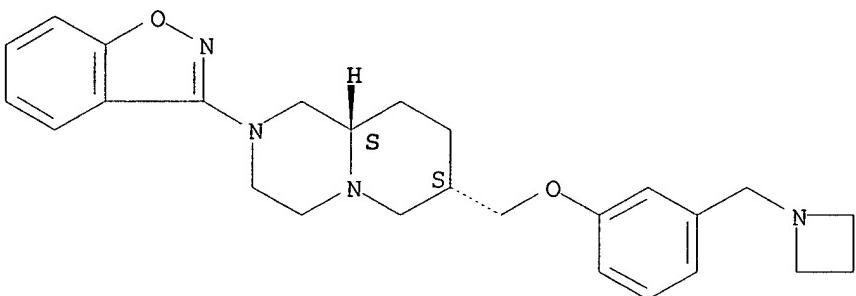
Absolute stereochemistry.



RN 247091-38-9 CAPLUS

CN 2H-Pyrido[1,2-a]pyrazine, 7-[[3-(1-azetidinylmethyl)phenoxy]methyl]-2-(1,2-benzisoxazol-3-yl)octahydro-, (7S,9aS)- (9CI) (CA INDEX NAME)

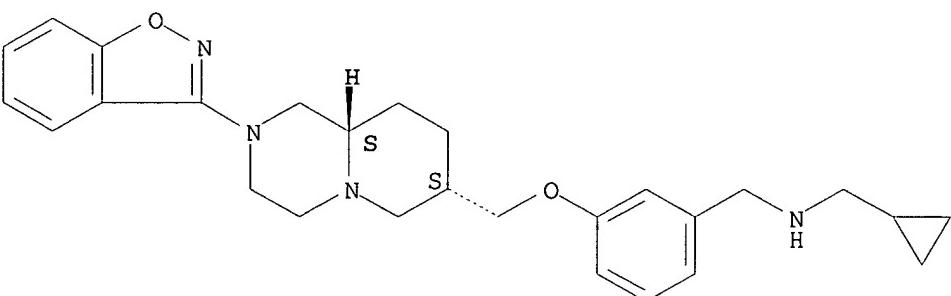
Absolute stereochemistry.



RN 247091-39-0 CAPLUS

CN Benzenemethanamine, 3-[[[(7S,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy]-N-(cyclopropylmethyl)- (9CI) (CA INDEX NAME)

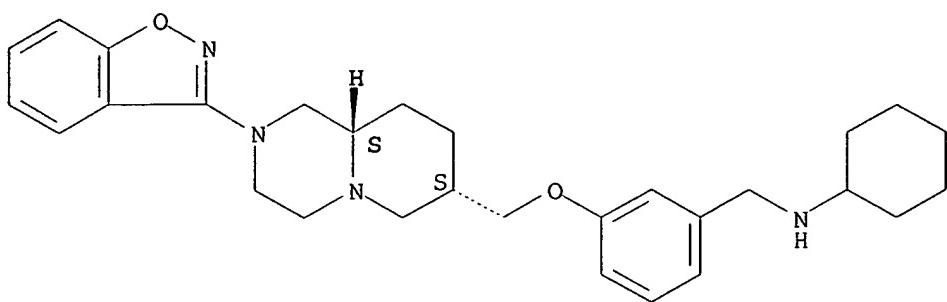
Absolute stereochemistry.



RN 247091-40-3 CAPLUS

CN Benzenemethanamine, 3-[[[(7S,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy]-N-cyclohexyl- (9CI) (CA INDEX NAME)

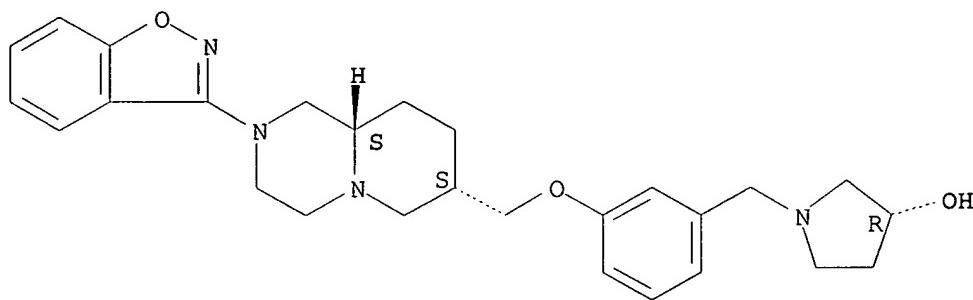
Absolute stereochemistry.



RN 247091-41-4 CAPLUS

CN 3-Pyrrolidinol, 1-[[3-[(7S,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy]phenyl]methyl]-, (3R)- (9CI) (CA INDEX NAME)

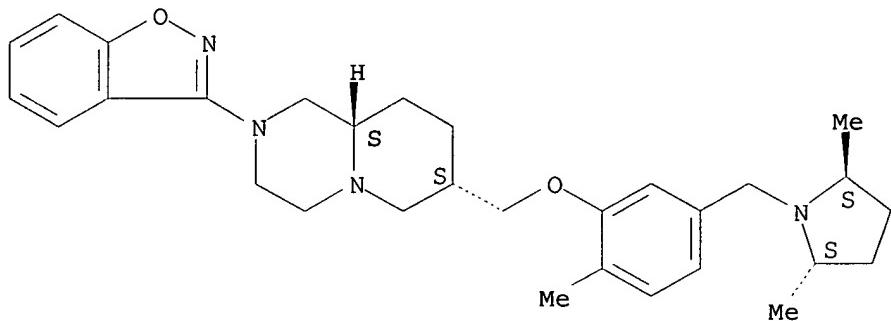
Absolute stereochemistry.



RN 247091-42-5 CAPLUS

CN 2H-Pyrido[1,2-a]pyrazine, 2-(1,2-benzisoxazol-3-yl)-7-[[5-[(2S,5S)-2,5-dimethyl-1-pyrrolidinyl]methyl]-2-methylphenoxy]methyl]octahydro-, (7S,9aS)- (9CI) (CA INDEX NAME)

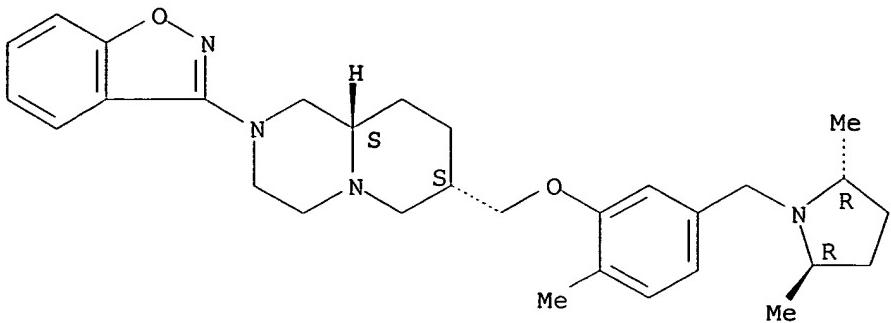
Absolute stereochemistry.



RN 247091-43-6 CAPLUS

CN 2H-Pyrido[1,2-a]pyrazine, 2-(1,2-benzisoxazol-3-yl)-7-[[5-[(2R,5R)-2,5-dimethyl-1-pyrrolidinyl]methyl]-2-methylphenoxy]methyl]octahydro-, (7S,9aS)- (9CI) (CA INDEX NAME)

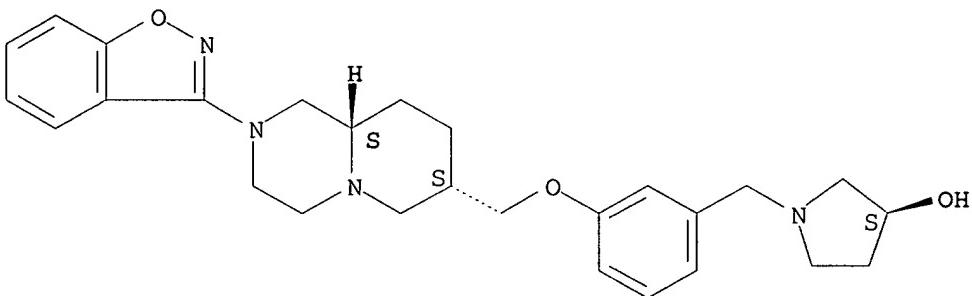
Absolute stereochemistry.



RN 247091-44-7 CAPLUS

CN 3-Pyrrolidinol, 1-[[3-[(7S,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy]phenyl]methyl]-, (3S)- (9CI) (CA INDEX NAME)

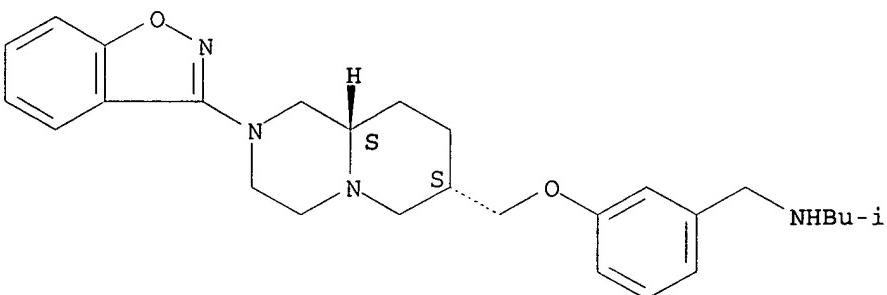
Absolute stereochemistry.



RN 247091-45-8 CAPLUS

CN Benzenemethanamine, 3-[[[(7S,9aS)-2-(1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy]-N-(2-methylpropyl)- (9CI) (CA INDEX NAME)

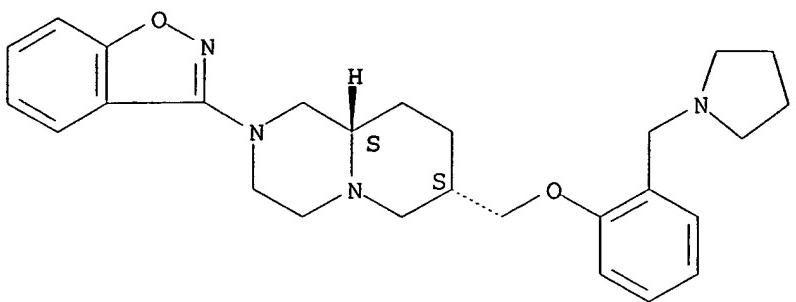
Absolute stereochemistry.



RN 247091-47-0 CAPLUS

CN 2H-Pyrido[1,2-a]pyrazine, 2-(1,2-benzisoxazol-3-yl)octahydro-7-[[2-(1-pyrrolidinylmethyl)phenoxy]methyl]-, (7S,9aS)- (9CI) (CA INDEX NAME)

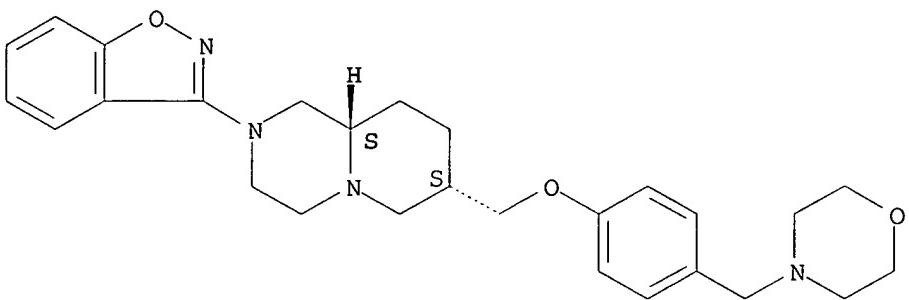
Absolute stereochemistry.



RN 247091-48-1 CAPLUS

CN 2H-Pyrido[1,2-a]pyrazine, 2-(1,2-benzisoxazol-3-yl)octahydro-7-[[4-(4-morpholinylmethyl)phenoxy]methyl]-, (7S,9aS)- (9CI) (CA INDEX NAME)

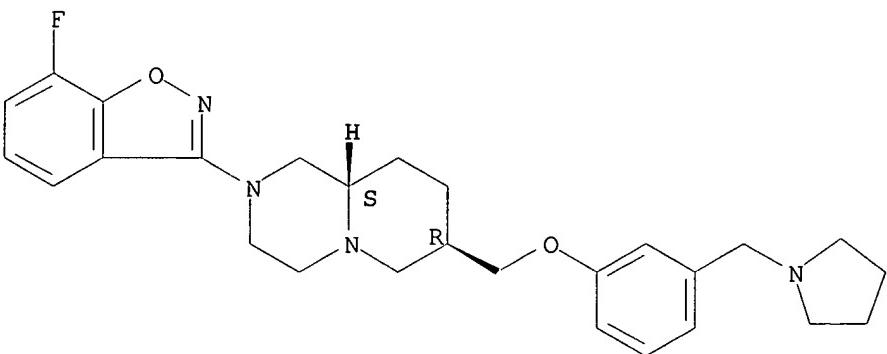
Absolute stereochemistry.



RN 247091-49-2 CAPLUS

CN 2H-Pyrido[1,2-a]pyrazine, 2-(7-fluoro-1,2-benzisoxazol-3-yl)octahydro-7-[[3-(1-pyrrolidinylmethyl)phenoxy]methyl]-, (7R,9aS)- (9CI) (CA INDEX NAME)

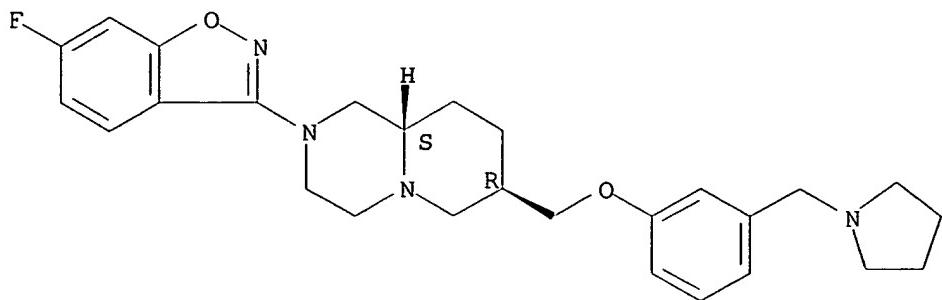
Absolute stereochemistry.



RN 247091-50-5 CAPLUS

CN 2H-Pyrido[1,2-a]pyrazine, 2-(6-fluoro-1,2-benzisoxazol-3-yl)octahydro-7-[[3-(1-pyrrolidinylmethyl)phenoxy]methyl]-, (7R,9aS)- (9CI) (CA INDEX NAME)

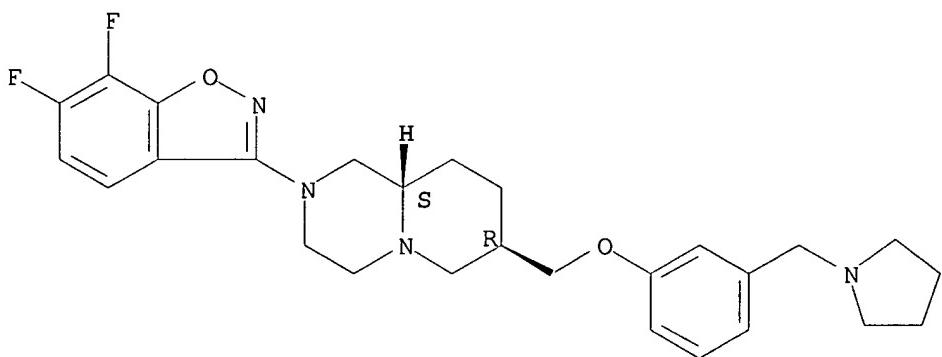
Absolute stereochemistry.



RN 247091-51-6 CAPLUS

CN 2H-Pyrido[1,2-a]pyrazine, 2-(6,7-difluoro-1,2-benzisoxazol-3-yl)octahydro-7-[(3-(1-pyrrolidinylmethyl)phenoxy)methyl]-, (7R,9aS)- (9CI) (CA INDEX NAME)

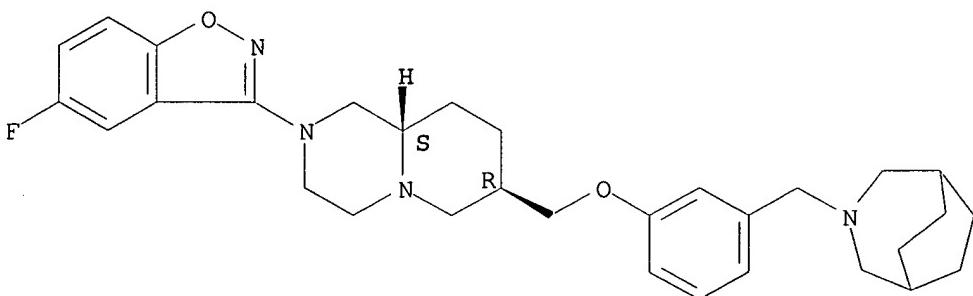
Absolute stereochemistry.



RN 247091-52-7 CAPLUS

CN 3-Azabicyclo[3.2.2]nonane, 3-[[3-[(7R,9aS)-2-(5-fluoro-1,2-benzisoxazol-3-yl)octahydro-2H-pyrido[1,2-a]pyrazin-7-yl]methoxy]phenyl]methyl]- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

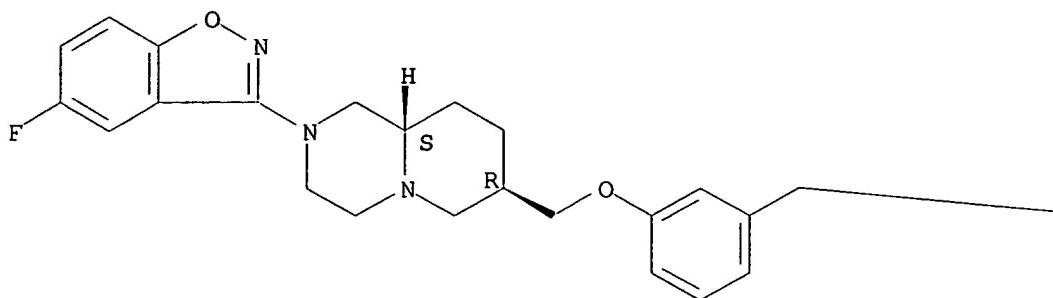


RN 247091-53-8 CAPLUS

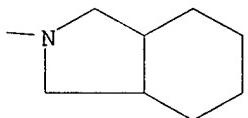
CN 2H-Pyrido[1,2-a]pyrazine, 2-(5-fluoro-1,2-benzisoxazol-3-yl)octahydro-7-[(3-[(octahydro-2H-isoindol-2-yl)methyl]phenoxy)methyl]-, (7R,9aS)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

PAGE 1-A



PAGE 1-B



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LOGOFF? (Y)/N/HOLD:Y

COST IN U.S. DOLLARS

SINCE FILE ENTRY	TOTAL SESSION
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13.35	180.50
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FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE ENTRY	TOTAL SESSION
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-2.25	-2.25
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